## IN THE CLAIMS:

This listing of claims will replace all prior versions, and listings, of claims in the application:

## **Listing of Claims:**

## 1. (Currently Amended) A compound of formula I

## wherein

- R<sub>1</sub> and R<sub>4</sub> are each independently H, C<sub>1</sub>-C<sub>10</sub>alkyl optionally substituted with one or more halogen, hydroxy, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>6</sub>, CONR<sub>7</sub>R<sub>8</sub>, C<sub>3</sub>-C<sub>7</sub>cycloalkyl or optionally substituted phenyl groups, or
  - phenyl optionally substituted with one to three halogen, hydroxy, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>9</sub>, NR<sub>10</sub>R<sub>11</sub> or CN groups;
- R<sub>2</sub> is H, C<sub>1</sub>-C<sub>6</sub>alkyl optionally substituted with a phenyl, naphthyl or <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S</u> group each group optionally substituted with one to three C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, hydroxy, CHO, NO<sub>2</sub>, CN, CO<sub>2</sub>R<sub>12</sub> or NR<sub>13</sub>R<sub>14</sub> groups,
  - phenyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>- $C_6$ alkyl, C<sub>1</sub>- $C_6$ haloalkyl, C<sub>1</sub>- $C_6$ alkoxy, phenyl, phenoxy, benzyl, benzyloxy,  $CO_2R_{17}$ ,  $NR_{18}R_{19}$  or  $CH_2CO_2R_{20}$  groups,
  - naphthyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, CO<sub>2</sub>R<sub>17</sub>, NR<sub>18</sub>R<sub>19</sub> or CH<sub>2</sub>CO<sub>2</sub>R<sub>20</sub> groups,
  - C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>17</sub> or NR<sub>18</sub>R<sub>19</sub> groups, or 5- to 10-membered heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms</u> selected from N, O or S optionally substituted with one to three halogen,

NO<sub>2</sub>, CN, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>17</sub> or NR<sub>18</sub>R<sub>19</sub> groups;

- R<sub>3</sub> is phenyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups,
  - $\underline{C_5}$ - $\underline{C_7}$ cycloheteroalkyl optionally substituted with one or more halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>,R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups, or
  - 5- to 10-membered heteroaryl ring system containing 1, 2 or 3 heteroatoms selected from N, O or S optionally substituted with one or more halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups;
  - R<sub>6</sub>, R<sub>9</sub>, R<sub>12</sub>, R<sub>17</sub>, R<sub>20</sub> and R<sub>26</sub> are each independently H or a C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, phenyl, C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl or <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S group each optionally substituted;</u>

n is 0 or an integer of 1 or 2; and

R<sub>7</sub>, R<sub>8</sub>, R<sub>10</sub>, R<sub>11</sub>, R<sub>13</sub>, R<sub>14</sub>, R<sub>18</sub>, R<sub>19</sub>, R<sub>21</sub>, R<sub>22</sub>, R<sub>24</sub> and R<sub>25</sub> are each independently H or a C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>3</sub>-C<sub>7</sub>cycloalkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, phenyl, C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl or <u>5</u>-to <u>10</u>-membered heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S</u> group each optionally substituted or each of R<sub>7</sub> and R<sub>8</sub> or R<sub>10</sub> and R<sub>11</sub> or R<sub>13</sub> and R<sub>14</sub> or R<sub>18</sub> and R<sub>19</sub> or R<sub>21</sub> and R<sub>22</sub> or R<sub>24</sub> and R<sub>25</sub> may be taken together with the nitrogen atom to which they are attached to form a 5- to 7-membered ring optionally containing another heteroatom selected from O, N or S; or

the stereoisemers thereof or the pharmaceutically acceptable salts thereof.

- 2. (Currently amended) The compound according to claim 1 wherein  $R_2$  is an optionally substituted phenyl or <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S</u> group.
- 3. (Original) The compound according to claim 1 wherein R<sub>1</sub> is H, C<sub>1</sub>-C<sub>3</sub>alkyl or an optionally substituted benzyl group.
- 4. (Currently Amended) The compound according to claim 1 wherein  $R_3$  is a  $C_5$ - $C_7$ cycloheteroalkyl, <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3</u>

heteroatoms selected from N, O or S or phenyl group each optionally substituted with one or two halogen, CN, NO<sub>2</sub>, CF<sub>3</sub>, methoxy, carboxy or SOR<sub>26</sub> groups.

- 5. (Original) The compound according to claim 2 wherein R<sub>4</sub> is H or phenyl or C<sub>1</sub>-C<sub>4</sub>alkyl optionally substituted with one hydroxy or phenyl group.
- 6. (Original) The compound according to claim 2 wherein  $R_3$  is a thienyl, pyridyl or phenyl group, each optionally substituted with one or two halogen, CN,  $NO_2$ ,  $CF_3$ , methoxy, carboxy or SOCH<sub>3</sub> groups.
- 7. (Original) The compound according to claim 3 wherein  $R_2$  is a phenyl group substituted with one or two halogen.
- 8. (Original) The compound according to claim 7 wherein  $R_3$  is a phenyl group substituted with one  $NO_2$  or  $CF_3$  group.
- 9. (Original) The compound according to claim 8 wherein  $R_1$  is H and  $R_4$  is H or  $CH_3$ .
  - 10. (Cancelled)
  - 11. (Cancelled)
  - 12. (Cancelled)
  - 13. (Cancelled)
  - 14. (Cancelled)
- 15. (Currently Amended) A pharmaceutical composition which comprises a pharmaceutically acceptable carrier and an effective amount of a compound of formula I

**(I)** 

wherein

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R<sub>1</sub> and R<sub>4</sub> are each independently H, C<sub>1</sub>-C<sub>10</sub>alkyl optionally substituted with one or more halogen, hydroxy, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>6</sub>, CONR<sub>7</sub>R<sub>8</sub>, C<sub>3</sub>-C<sub>7</sub>cycloalkyl or optionally substituted phenyl groups, or

- phenyl optionally substituted with one to three halogen, hydroxy, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>9</sub>, NR<sub>10</sub>R<sub>11</sub> or CN groups;
- R<sub>2</sub> is H, C<sub>1</sub>-C<sub>6</sub>alkyl optionally substituted with a phenyl, naphthyl or <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S</u> group each group optionally substituted with one to three C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, hydroxy, CHO, NO<sub>2</sub>, CN, CO<sub>2</sub>R<sub>12</sub> or NR<sub>13</sub>R<sub>14</sub> groups,
  - phenyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>- $C_6$ alkyl, C<sub>1</sub>- $C_6$ haloalkyl, C<sub>1</sub>- $C_6$ alkoxy, phenyl, phenoxy, benzyl, benzyloxy, CO<sub>2</sub>R<sub>17</sub>, NR<sub>18</sub>R<sub>19</sub> or CH<sub>2</sub>CO<sub>2</sub>R<sub>20</sub> groups,
  - naphthyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, CO<sub>2</sub>R<sub>17</sub>, NR<sub>18</sub>R<sub>19</sub> or CH<sub>2</sub>CO<sub>2</sub>R<sub>20</sub> groups,
  - C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>17</sub> or NR<sub>18</sub>R<sub>19</sub> groups, or 5- to 10-membered heteroaryl ring system containing 1, 2 or 3 heteroatoms selected from N, O or S optionally substituted with one to three halogen, NO<sub>2</sub>, CN, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>4</sub>alkoxy, CO<sub>2</sub>R<sub>17</sub> or NR<sub>18</sub>R<sub>19</sub> groups;
- R<sub>3</sub> is phenyl optionally substituted with one to three halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups,
  - C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl optionally substituted with one or more halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>,R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups, or
  - <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms</u> selected from N, O or S optionally substituted with one or more halogen, NO<sub>2</sub>, CN, hydroxy, C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>1</sub>-C<sub>6</sub>alkylthio, C<sub>1</sub>-C<sub>6</sub>haloalkyl, C<sub>1</sub>-C<sub>6</sub>alkoxy, phenyl, phenoxy, benzyl, benzyloxy, SO<sub>n</sub>R<sub>26</sub>, SO<sub>2</sub>NR<sub>21</sub>R<sub>22</sub>, CO<sub>2</sub>R<sub>23</sub> or NR<sub>24</sub>R<sub>25</sub> groups;
  - R<sub>6</sub>, R<sub>9</sub>, R<sub>12</sub>, R<sub>17</sub>, R<sub>20</sub> and R<sub>26</sub> are each independently H or a C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, phenyl, C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl or <u>5- to 10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S group each optionally substituted;</u>

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n is 0 or an integer of 1 or 2; and

R<sub>7</sub>, R<sub>8</sub>, R<sub>10</sub>, R<sub>11</sub>, R<sub>13</sub>, R<sub>14</sub>, R<sub>18</sub>, R<sub>19</sub>, R<sub>21</sub>, R<sub>22</sub>, R<sub>24</sub> and R<sub>25</sub> are each independently H or a C<sub>1</sub>-C<sub>6</sub>alkyl, C<sub>3</sub>-C<sub>7</sub>cycloalkyl, C<sub>1</sub>-C<sub>6</sub>haloalkyl, phenyl, C<sub>5</sub>-C<sub>7</sub>cycloheteroalkyl or <u>5</u>-to <u>10-membered</u> heteroaryl <u>ring system containing 1, 2 or 3 heteroatoms selected from N, O or S</u> group each optionally substituted or each of R<sub>7</sub> and R<sub>8</sub> or R<sub>10</sub> and R<sub>11</sub> or R<sub>13</sub> and R<sub>14</sub> or R<sub>18</sub> and R<sub>19</sub> or R<sub>21</sub> and R<sub>22</sub> or R<sub>24</sub> and R<sub>25</sub> may be taken together with the nitrogen atom to which they are attached to form a 5- to 7-membered ring optionally containing another heteroatom selected from O, N or S; or

the stereoisomers thereof or the pharmaceutically acceptable salts thereof.

- 16. (Original) The composition according to claim 15 having a formula I compound wherein R<sub>2</sub> is an optionally substituted phenyl, thienyl or pyridyl group.
- 17. (Original) The composition according to claim 16 having a formula I compound wherein  $R_1$  is H and  $R_4$  is H or CH<sub>3</sub>.
- 18. (Original) The composition according to claim 17 having a formula I compound wherein R<sub>3</sub> is a thienyl, pyridyl or phenyl group each optionally substituted with one or two halogen, CN, NO<sub>2</sub>, CF<sub>3</sub>, methoxy, carboxy or SOCH<sub>3</sub> groups.
- 19. (Original) The composition according to claim 18 having a formula I compound wherein R<sub>2</sub> is a phenyl group substituted with one or two halogen.
- 20. (Currently Amended) The composition according to claim 15 having a formula I compound selected from the group consisting of:
- 2-(3-fluorophenyl)-4-(3-nitrophenyl)-1,6-dihydrodipyrazolo[3,4-b:3',4'-d]pyridin-3(2H)-one;
- 2-(3-fluorophenyl-6-methyl-4-(3-nitrophenyl)-1,6-dihydrodipyrazolo[3,4-b:3',4'-d]-pyridin-3(2H)-one;
- 2-(4-chlorophenyl)-6-methyl-4-[3-(trifluoromethyl)phenyl]-1,6-dihydrodipyrazolo-[3,4-b:3',4'-d]pyridin-3(2H)-one;
- 2-(4-chlorophenyl)-6-methyl-4-(3-fluorophenyl)-1,6-dihydrodipyrazolo-[3,4-b:3'4'-d]pyridin-3(2H)-one;
- 4-(5-bromo-3-pyridinyl)-6-methyl-3-[(trifluoromethyl)phenyl]-1,6-dihydrodipyrazolo-[3,4-b:3',4'-d]pyridin-3(2H)-one;
- 4-(5-bromo-3-pyridinyl)-3-(4-fluorophenyl)-6-methyl-1,6-dihydrodipyrazolo-[3,4-b:3',4'-d]pyridin-3-(2H)-one;
- methyl 3-{6-methyl-3-oxo-4-[3-(trifluoromethyl)phenyl]-3,6-dihydrodipyrazolo-[3,4-b:3',4'-d]pyridin-2(1H)-yl}benzoate;
- 2-chloro-5-{6-methyl-3-oxo-4-[3-(trifluoromethyl)phenyl]-3,6-dihydrodipyrazolo-

[3,4-b:3',4'-d]pyridin-2(1H)-yl}benzoic acid;

- 4-(3-bromophenyl)-6-methyl-2-(4-nitrophenyl)-1,6-dihydrodipyrazolo[3,4-b:3',4'-d]-pyridin-3(2H)-one;
- 4-[4-(3-bromophenyl)-6-methyl-3-oxo-3,6-dihdrodipyrazolo[3,4-b:3',4'-d]pyridin-2(1H)-yl]-2-chlorobenzoic acid;
- methyl 2-fluoro-4-{6-methyl-3-oxo-4-[3-(trifluoromethyl)phenyl]-3,6-dihydrodipyrazolo-[3,4-b:3', 4'-d]pyridin-2-(1H)-yl}benzoate;

the stereoisomers thereof;

and the pharmaceutically acceptable salts thereof.